L Number	Hits	Search Text	DB	Time stamp
1	568	lysine same polyamide	USPAT;	2004/10/20 15:34
-		7,55 52 p 5.7,4	US-PGPUB;	200 1/10/20 13.54
			EPO; JPO;	
		•	DERWENT	
2	484	(lysine same polyamide) and polymer		2004/10/20 15:25
2	707		USPAT;	2004/10/20 15:35
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
3	1	((lysine same polyamide) and polymer) and geminal	USPAT;	2004/10/20 15:38
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
4	176	((lysine same polyamide) and polymer) and dimer	USPAT;	2004/10/20 15:39
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
5	0	PCT/JP03/05453	USPAT;	2004/10/20 15:39
			US-PGPUB;	
		•	EPO; JPO;	
			DERWENT	
6	o	"PCT/JP03/05453"	USPAT;	2004/10/20 15:39
			US-PGPUB;	200 1/10/20 13.39
			EPO; JPO;	
			DERWENT	
7	2196	hanabusa.in.		2004/10/20 15:20
'	2190	Hallabusa.iii.	USPAT;	2004/10/20 15:39
		T.	US-PGPUB;	
			EPO; JPO;	
	25222	· · · · · · · · · · · · · · · · · · ·	DERWENT	7004/40/20 45 20
8	252333	suzuki.in.	USPAT;	2004/10/20 15:39
			US-PGPUB;	
			EPO; JPO;	
		l	DERWENT	
9	41	hanabusa.in. and suzuki.in.	USPAT;	2004/10/20 15:42
			US-PGPUB;	
	ļ		EPO; JPO;	
			DERWENT	
10	6261	514/2.ccls.	USPAT;	2004/10/20 15:44
			US-PGPUB;	
			EPO; JPO;	
			DERWENT	
11	3308	530/300.ccls.	USPAT;	2004/10/20 15:48
			US-PGPUB;	
į			EPO; JPO;	
			DERWENT	
12	675	530/323,332.ccls.	USPAT;	2004/10/20 15:49
		, ,	US-PGPUB;	
			EPO; JPO;	
			DERWENT	
13	2283	554/112,106,69,66,57,56,47.ccls. 564/153,152.ccls.	USPAT;	2004/10/20 15:49
20	2203		US-PGPUB;	200 1/10/20 13.43
			EPO; JPO;	' '
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14	6182	530/300.ccls. 530/323,332.ccls.		2004/10/20 15:50
177	0102	550/300.ccis. 550/323,332.ccis. (554/112,106,69,66,57,56,47.ccis. 564/153,152.ccis.)	USPAT;	2004/10/20 15:50
		(33,152,00,03,00,37,30,47,00,5,304/153,152,00)	US-PGPUB;	
			EPO; JPO;	
15	272	(F30/300 celo, F30/333 333!-	DERWENT	2004//2/20 := ==
15	273	(530/300.ccls. 530/323,332.ccls.	USPAT;	2004/10/20 15:50
		(554/112,106,69,66,57,56,47.ccls. 564/153,152.ccls.)) and	US-PGPUB;	
		chelator	EPO; JPO;	
		·	DERWENT	

1.0		(/===/===		
16	140	((530/300.ccls. 530/323,332.ccls.	USPAT;	2004/10/20 15:50
		(554/112,106,69,66,57,56,47.ccls. 564/153,152.ccls.)) and	US-PGPUB;	
		chelator) and polymer	EPO; JPO;	
			DERWENT	
17	127	(((530/300.ccls. 530/323,332.ccls.	USPAT;	2004/10/20 15:50
		(554/112,106,69,66,57,56,47.ccls. 564/153,152.ccls.)) and	US-PGPUB;	
		chelator) and polymer) and gel	EPO; JPO;	
			DERWENT	
18	57	((((530/300.ccls. 530/323,332.ccls.	USPAT;	2004/10/20 15:50
		(554/112,106,69,66,57,56,47.ccls. 564/153,152.ccls.)) and	US-PGPUB;	
		chelator) and polymer) and gel) and dimer	EPO; JPO;	
			DERWENT	

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Enter NEWS followed by the item number or name to see news on that

CAS World Wide Web Site (general information)

35 15 26 -32 29 Ĥ 30 27 38

29 26-33 29-32 32-35 32-37 33-34 33-36 exact bonds:
1-38 2-4 4-5 4-19 4-20 7-39 8-10 8-12 8-18 9-11 9-13 9-17 17-24 17-25 18-26 18-27 28-31 29-30

Match level:
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS
8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS
15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 23:CLASS
24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS
31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS 36:CLASS 37:CLASS
38:CLASS 39:CLASS

## L1 STRUCTURE UPLOADED

SAMPLE SEARCH INITIATED 15:17:49 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 16171 TO ITERATE

6.2% PROCESSED 1000 ITERATIONS ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

ONLINE \*\*COMPLETE\*\*
BATCH \*\*COMPLETE\*\*
315807 TO 331033
0 TO 0 FULL FILE PROJECTIONS: PROJECTED ITERATIONS: PROJECTED ANSWERS:

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=> fil reg COST IN U.S. DOLLARS TOTAL SINCE FILE FNTRY FULL ESTIMATED COST 0.21 0.21

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STRUCTURE FILE UPDATES: 19 OCT 2004 HIGHEST RN 765878-56-6 DICTIONARY FILE UPDATES: 19 OCT 2004 HIGHEST RN 765878-56-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

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Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: <a href="http://www.cas.org/ONLINE/DBSS/registryss.html">http://www.cas.org/ONLINE/DBSS/registryss.html</a>

Uploading H:\STN queries\10777179.str

· 0 SEA SSS SAM L1

s 11 ful FULL SEARCH INITIATED 15:17:53 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 321251 TO ITERATE

87.4% PROCESSED 280616 ITERATIONS ANSWERS

100.0% PROCESSED 321251 ITERATIONS

21

21

SEARCH TIME: 00.00.26

21 SEA SSS FUL L1

=> file hcaplus COST IN U.S. DOLLARS TOTAL SINCE FILE ENTRY SESSION FULL ESTIMATED COST 155.84

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FILE COVERS 1907 - 20 Oct 2004 VOL 141 ISS 1 FILE LAST UPDATED: 19 Oct 2004 (20041019/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.  $\label{eq:contains} % \begin{subarray}{ll} \end{subarray} % \begin{subarray}{ll} \end{subarra$ 

=> s 13 ∟4 4 L3

=> d 14 1-4 ibib abs ANSWER 1 OF 4 HCAPLUS COPYRIGHT 2004 ACS ON STN SSION NUMBER: 2003:878000 HCAPLUS Full-text 140:181736 ACCESSION NUMBER: DOCUMENT NUMBER:

specific topic.

TITLE: L-Lysine based gemini organogelators: their organogelation properties and thermally stable organogels AUTHOR(S): Sužuki, Masahiro; Nigawara, Tomomi; Yumoto, Kimura, Mutsumi; Shirai, Hirofusa; Hanabusa, Kenji CORPORATE SOURCE: Shinshu Graduate School of Science and Technology, University, Ueda, Nagano, 386-8567, Japa Organic & Biomolecular Chemistry (2003) SOURCE: CODEN: OBCRAK; ISSN: 1477-0520 Royal Society of Chemistry Journal Royal Society of Chemistry

DOCUMENT TYPE: Journal
LANGUAGE: English

AB Novel gemini organogelators based on L-lysine, in which two Llysine derivs. are linked by different alkylene chain lengths
through the amide bond, have been simply and effectively
synthesized, and their organogelation abilities and thermal
stabilities have been investigated. In a series of L-lysine Et
ester derivs., the organogelation abilities decreased with
increasing alkylene spacer length. In particular, bis(Nelauroyl-L-lysine Et ester)oxalyl amide,
H23C11CONH(CH2)4CH(COZET)NH-COCO-NHCH(COZET)(CH2)4NHCOC11H23, is
a good organogelator that gels most organic solvents such as
alcs., cyclic ethers, aromatic solvents and acetonitrile.
Various oxalyl amide derivs. with different alkyl ester groups
such as hexyl decyl, dodecyl, 2-ethyl-l-hexyl and 3,5,5trimethylhexyl also showed good organogelation abilities.
Furthermore, it was found that the cyclohexane gels formed by
some oxalyl amide derivs. have a high thermal stability.
REFERENCE COUNT: 40 THERE ARE 40 CITED REFERENCES

RECORD. ALL CITATIONS AVAILABLE THERE

REFORMAT PUBLISHER:

L4 ANSWER 2 OF 4 HCAPLUS COPYRIGHT 2004 ACS ON STN ACCESSION NUMBER: 2003:627026 HCAPLUS Full-text 139:337687 TITLE: New gowie:

New gemini organogelators linked by oxalyl amide:

organogel formation and their thermal

stabilities AUTHOR(S): Mariko;

Suzuki, Masahiro; Nigawara, Tomomi; Yumoto,

Kimura, Mutsumi; Shirai, Hirofusa; Hanabusa,

Kenji CORPORATE SOURCE: Graduate School of Science and Technology,

Shinshu SOURCE: 6843

University, Ueda, Nagano, 386-8567, Japan Tetrahedron Letters (2003), 44(36), 6841-

CODEN: TELEAY; ISSN: 0040-4039

Optically active polyamides with regular structural sequences were prepd. from L-lysine and adipic acid. An optically active sym. diamine, N,N'-bis(L-5-amino-5-carboxyamyl) adipamide, m. 305' (decomposition), was obtained by treating L-lysine with adipoyl chloride (1) in the presence of Cu2+. The interfacial polycondensation of this diamine with I gave a regular polymer, while the polycondensation of L-lysine with I gave an irregular polymer of the same anal. composition The m.ps., optical rotations, and the O.R.D. curves of these polymers were investigated. ΑB investigated.

=> FIL STNGUIDE COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 168.61 12.56 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) TOTAL SINCE FILE **ENTRY** SESSION -2.80 ့ CA SUBSCRIBER PRICE 2.80

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FILE CONTAINS CURRENT INFORMATION. LAST RELOADED: Oct 15, 2004 (20041015/UP).

=> FIL REGISTRY COST IN U.S. DOLLARS

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PUBLISHER: DOCUMENT TYPE: LANGUAGE:

SHER: Elsevier Science B.V.

IENT TYPE: Journal

MGE: English

SOURCE(S): CASREACT 139:337687

New gemini organogelators linked by an oxalyl amide that can be easily, effectively, and cheaply synthesized have good organogelation abilities and their cyclohexane gels have superior thermal stabilities; especially 7 possessing the branched alkyl ester can gel at 0.7 wt% cyclohexane even at OTHER SOURCE(S):

70°C. REFERENCE COUNT: AVAILABLE FOR THIS

36 THERE ARE 36 CITED REFERENCES

RECORD, ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L4 ANSWER 3 OF 4 HCAPLUS COPYRIGHT 2004 ACS ON STN ACCESSION NUMBER: 1997:334995 HCAPLUS Full-text 127:51064 DOCUMENT NUMBER: TITLE:

Synthesis and characterization of random and regular

AUTHOR(S): Sekiguchi, Hikaru CORPORAȚE SOURCE:

L-lysine-based polyamides Gachard, Isabelle; Coutin, Bernard;

Laboratoire Chimie Macromoleculaire,

SOURCE: 198(5),

Pierre et Marie Curie, Paris, F-75252, Fr. Macromolecular Chemistry and Physics (1997),

1375-1389

PUBLISHER: DOCUMENT TYPE: LANGUAGE:

1375-1389
CODEN: MCHPES; ISSN: 1022-1352
Huethig & Wepf
Journal
JAGE: English
The synthesis of polyamides based on the natural diamine Llysine and diacids, adipic or glutaric acid, is described. They
were obtained by polycondensation of active diesters,
pentachlorophenyl, and pentafluorophenyl esters. L-Lysine being
non-sym., aregular (random), and syndioregular (head-to-head,
tail-to-tail) poly(adipoyl-L-lysine)s and poly(glutaroyl-Llysine)s were obtained with mol. wts. > 15,000 while isoregular
(head-to-tail) poly(adipoyl-L-lysine)s and poly(glutaroyl-Llysine)s were prepared with lower mol. wts.

L4 ANSWER 4 OF 4 HCAPLUS COPYRIGHT 2004 ACS on STN ACCESSION NUMBER: 1968:13536 HCAPLUS Full-text 68:13536 Optically active polyamides with regular

structural AUTHOR(S): CORPORATE SOURCE:

SOURCE:

sequences prepared from α-amino acids Saotome, Kazuo; Schulz, Rolf Christian Univ., Mainz, Mainz, Fed. Rep. Ger. Makromolekulare Chemie (1967), 109, 239-48 CODEN: MACEAK; ISSN: 0025-116X Journal

DOCUMENT TYPE: LANGUAGE:

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

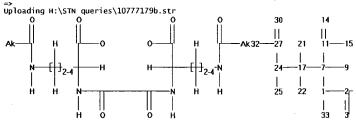
STRUCTURE FILE UPDATES: 19 OCT 2004 HIGHEST RN 765878-56-6 DICTIONARY FILE UPDATES: 19 OCT 2004 HIGHEST RN 765878-56-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting  ${\sf SmartSELECT}$  searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: <a href="http://www.cas.org/ONLINE/DBSS/registryss.html">http://www.cas.org/ONLINE/DBSS/registryss.html</a>



chain nodes: 1 2 3 4 5 21 22 23 24  $\begin{smallmatrix} 6 & 7 & 8 & 9 & 10 & 11 & 12 & 13 & 14 & 15 & 16 & 17 \\ 25 & 26 & 27 & 28 & 29 & 30 & 31 & 32 & 33 & 34 \end{smallmatrix}$ 18 19 20 17-22 17-24 23-26 23-26 24-27 24-27 27-30 27-32 26-29 26-31 exact/norm bonds :
1-2 1-7 2-3 4-6 4-5 6-8 11-14 11-15 12-13 12-18 16-23 17-24 23-28 24-27 27-30 27-32 28-29 28-31 24 25-28 24-27 27-30 27-32 28-29 28-31 exact bonds:
1-33 2-4 6-34 7-9 7-11 7-17 8-10 8-12 8-16 16-19 16-20 17-21 17-22 23-26 24-25

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS
8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS

=> s 15 ful FULL SEARCH INITIATED 15:22:16 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 7829 TO ITERATE

100.0% PROCESSED 7829 ITERATIONS

ANSWERS SEARCH TIME: 00.00.01

7 SEA SSS FUL LS

=> file hcaplus COST IN U.S. DOLLARS TOTAL SINCE FILE ENTRY SESSION FULL ESTIMATED COST 324.69 155.84

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FILE COVERS 1907 - 20 Oct 2004 VOL 141 ISS 17 FILE LAST UPDATED: 19 Oct 2004 (20041019/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 16

RN 615584-81-1 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-82-2 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediy1)bis[N6-(1-oxododecy1)-, dihexy1 ester.(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-83-3 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, didecyl ester (9CI) (CA INDEX NAME)

2 L6

=> d 17 1-2 ibib abs hitstr

ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2004 ACS ON STN
SSION NUMBER: 2003:878000 HCAPLUS <u>Full-text</u>
MENT NUMBER: 140:181736
E: L-Lysine based gemini organogelators: their
organogelation properties and thermally DOCUMENT NUMBER: TITLE:

stable

organogels Suzuki, Masahiro; Nigawara, Tomomi; Yumoto, AUTHOR(S): Mariko;

Kimura, Mutsumi; Shirai, Hirofusa; Hanabusa, Graduate School of Science and Technology,

Kenji CORPORATE SOURCE: Shinshu

University, Ueda, Nagano, 386-8567, Japan Organic & Biomolecular Chemistry (2003),

PUBLISHER: DOCUMENT TYPE: LANGUAGE: AB NOVE 1 gen

4124-4131
CODEN: OBCRAK; ISSN: 1477-0520
ISHER: Royal Society of Chemistry
RENT TYPE: Journal
JAGE: English
Novel gemini organogelators based on L-lysine, in which two Llysine derivs. are linked by different alkylene chain lengths
through the amide bond, have been simply and effectively
synthesized, and their organogelation abilities and thermal
stabilities have been investigated. In a series of L-lysine Et
ester derivs., the organogelation abilities decreased with
increasing alkylene spacer length. In particular, bis(Neester derivs., the organogelation abilities decreased with increasing alkylene spacer length. In particular, bis(Na-lauroyl-L-lysine Et ester)oxalyl amide, H23C11CONH(CH2)4CH(CO2Et)NH-COCO-NHCH(CO2Et)(CH2)4NHCOC11H23, is a good organogelator that gels most organic solvents such as alcs., cyclic ethers, aromatic solvents and acetonitrile. Various oxalyl amide derivs. with different alkyl ester groups such as hexyl, decyl, dodecyl, 2-ethyl-1-hexyl and 3,5,5-trimethylhexyl also showed good organogelation abilities. Furthermore, it was found that the cyclohexane gels formed by some oxalyl amide derivs. have a high thermal stability. 615584-80-0P 615584-81-1P 615584-82-2P 615584-83-3P 615584-84-4P 615584-85-5P 615584-86-6P RL: PRP (Properties); SPN (Synthetic preparation); PREP

615584-86-6P
RL: PRP (Properties); SPN (Synthetic preparation); PREP
(Preparation)
(preparation, organogelation property and thermal stability
of bis-lysine
amides linked by alkylene chains)
RN 615584-80-0 HCAPLUS
CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1oxododecyl)- (9CI)
(CA INDEX NAME)

Absolute stereochemistry.

Absolute stereochemistry.

RN 615584-84-4 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, didodecyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-85-5 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(2-ethylhexyl) ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-86-6 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(3,5,5-trimethylhexyl) ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT: AVAILABLE FOR THIS

RE FORMAT

stabilities AUTHOR(S): Mariko;

Kenji CORPORATE SOURCE: Shinshu

SOURCE: 6843

PUBLISHER:

THERE ARE 40 CITED REFERENCES 40

RECORD. ALL CITATIONS AVAILABLE IN THE

L7 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2004 ACS on STN
ACCESSION NUMBER: 2003:627026 HCAPLUS Full-text
139:337687
ITILE: amide: - New gemin organogelators linked by oxalyl

organogel formation and their thermal

Suzuki, Masahiro; Nigawara, Tomomi; Yumoto, Kimura, Mutsumi; Shirai, Hirofusa; Hanabusa,

Graduate School of Science and Technology,

University, Ueda, Nagano, 386-8567, Japan Tetrahedron Letters (2003), 44(36), 6841-

CODEN: TELEAY; ISSN: 0040-4039 Elsevier Science B.V.

oxododecyl)-, dihexyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-83-3 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, didecyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-84-4 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, didodecyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

DOCUMENT TYPE: Journal
LANGUAGE: English
OTHER SOURCE(S): CASREACT 139:337687

AB New gemini organoge lators linked by an oxalyl amide that can be easily, effectively, and cheaply synthesized have good organoge lation abilities and their cyclohexane gels have superior thermal stabilities; especially 7 possessing the branched alkyl ester can gel at 0.7 wt% cyclohexane even at 70°C.

IT 615584-80-0P 615584-81-1P 615584-82-2P 615584-83-3P 615584-86-6P 615584-86-6P
BI: PRP (Properties): SPN (Synthetic preparation): PREP 615584-86-6P
RL: PRP (Properties); SPN (Synthetic preparation); PREP
(Preparation)
(NMR and FT-IR on gelation of prepared gemini oxalyl-amide organogelators)
RN 615584-80-0 HCAPLUS
CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododeyl)- (9C1)
(CA INDEX NAME)

Absolute stereochemistry.

RN 615584-81-1 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

615584-82-2 HCAPLUS L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-

RN 615584-85-5 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(2-ethylhexyl) ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-86-6 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(3,5,5-trimethylhexyl) ester (9CI) (CA INDEX NAME) Absolute stereochemistry.

REFERENCE COUNT: 36

THERE ARE 36 CITED REFERENCES

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29 28-33 29-32 32-35 33-34 exact bonds:

1-38 2-4 4-5 4-19 4-20 7-39 8-10 8-12 8-18 9-11 9-13 9-17 17-24 17-25 18-26 18-27 28-31 29-30 32-37 33-36 36-44 37-54 44-45 45-46 46-47 47-48 48-49 49-50 50-51 51-52 52-53 54-55 55-56 56-57 57-58 58-59 59-60 60-61 61-62 62-63

Match level:
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS
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15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 29:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 33:CLASS 34:CLASS 34:CLASS 35:CLASS 36:CLASS 36:CLASS

conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

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100.0% PROCESSED 342183 ITERATIONS

ANSWERS SEARCH TIME: 00.00.08

17 SEA SSS FUL L8

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## => d 110 1-2 ibib abs hitstr

L10 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2004 ACS on STN
ACCESSION NUMBER: 2003:878000 HCAPLUS Full-text
DOCUMENT NUMBER: 140:181736
L-Lysine based gemini organogelators: their organogelation properties and thermally

stable

AUTHOR(S): Mariko;

organogels Suzuki, Masahiro; Nigawara, Tomomi; Yumoto, Kimura, Mutsumi; Shirai, Hirofusa; Hanabusa,

Kenji CORPORATE SOURCE: Shinshu

Graduate School of Science and Technology,

University, Ueda, Nagano, 386-8567, Japan Organic & Biomolecular Chemistry (2003),

SOURCE: 1(22),

PUBLISHER:

DOCUMENT TYPE:

LANGUAGE:

AB Novel gemini organogelators based on L-lysine, in which two L-lysine derivs. are linked by different alkylene chain lengths through the amide bond, have been simply and effectively synthesized, and their organogelation abilities and thermal stabilities have been investigated. In a series of L-lysine Et ester derivs., the organogelation abilities decreased with increasing alkylene spacer length. In particular, bis(NE-lauroyl-L-lysine Et ester)oxalyl amide,

H23C11cONN(CH2)4CH(CO2ET)NH-COCO-NHCH(CO2ET)(CH2)4NHCOC11H23, is a good organogelator that gels most organic solvents such as alcs., cyclic ethers, aromatic solvents and acctonitrile.

Various oxalyl amide derivs. with different alkyl ester groups such as hexyl, decyl, dodecyl, 2-ethyl-1-hexyl and 3,5,5-trimethylhexyl also showed good organogelation abilities.

Furthermore, it was found that the cyclohexane gels formed by some oxalyl amide derivs. have a high thermal stability.

II 615584-80-0P 615584-81-1P 615584-82-2P 615584-85-66 658051-85-67 658051-85-7P 658051-85-67 658051-95-7P 658

RN 615584-83-3 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediy1)bis[N6-(1-oxododecy1)-, didecy1 ester (9C1) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-84-4 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, diodecyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-85-5 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododeyl)-, bis(2-ethylhexyl) ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

(CA INDEX NAME)

Absolute stereochemistry.

RN 615584-81-1 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-82-2 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, dihexyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-86-6 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(3,5,5-trimethylhexyl) ester (9CI) (CA INDEX NAME) Absolute stereochemistry.

RN 658051-84-4 HCAPLUS CN L-Lysine, N2,N2'-(1,3-dioxo-1,3-propanediy1)bis[N6-(1-oxododecy1)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 658051-85-5 HCAPLUS CN L-Lysine, N2,N2'-(1,4-dioxo-1,4-butanediyl)bis[N6-(1oxododecyl)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

\_ (CH2)10 Me

RN 658051-86-6 HCAPLUS CN L-Lysine, N2,N2'-(1,5-dioxo-1,5-pentanediy1)bis[N6-(1-oxododecy1)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

\_ (CH2)10 Me

RN 658051-93-5 HCAPLUS CN L-Lysine, N2,N2'-(1,3-dioxo-1,3-propanediyl)bis[N6-(1-oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

~ (CH2)10 Me

RN 658051-94-6 HCAPLUS CN L-Lysine, N2,N2'-(1,4-dioxo-1,4-butanediyl)bis[N6-(1-oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

\_ (CH2)10 Me

RN 658051-87-7 HCAPLUS CN L-Lysine, N2,N2'-(1,6-dioxo-1,6-hexanediy1)bis[N6-(1-oxododecy1)- (9CT) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

\_\_(CH2)10\_\_\_\_\_Me

RN 658051-88-8 HCAPLUS
CN L-Lysine, N2,N2'-(1,7-dioxo-1,7-heptanediy1)bis[N6-(1-oxododecy1)- (9CI)
(CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

─(CH2)10 Me

RN 658051-95-7 HCAPLUS CN L-Lysine, N2,N2'-(1,5-dioxo-1,5-pentanediyl)bis[N6-(1oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

 ${\bf Absolute\ stereochemistry.}$ 

PAGE 1-B

~(CH2)10 Ме

RN 658051-96-8 HCAPLUS
CN L-Lysine, N2.N2'-(1,6-dioxo-1,6-hexanediyl)bis[N6-(1-oxododecyl)- diethyl ester (9C1) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

\_ (СН2)10 Ме

RN 658051-97-9 HCAPLUS CN L-Lysine, N2,N2'-(1,7-dioxo-1,7-heptanediy1)bis[N6-(1-oxododecy1)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-B

∼(CH2)10 Me

REFERENCE COUNT: AVAILABLE FOR THIS

THERE ARE 40 CITED REFERENCES

RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

L10 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2004 ACS ON STN ACCESSION NUMBER: 2003:627026 HCAPLUS Full-text 139:337687

RN 615584-82-2 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, dihexyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Absolute stereochemistry.

TITLE: amide: New gemini organogelators linked by oxalyl organogel formation and their thermal stabilities Suzuki, Masahiro; Nigawara, Tomomi; Yumoto. Kimura, Mutsumi; Shirai, Hirofusa; Hanabusa. Kenji CORPORATE SOURCE: Shinshu Graduate School of Science and Technology, University, Ueda, Nagano, 386-8567, Japan Tetrahedron Letters (2003), 44(36), 6841-SOURCE:
6843

CODEN: TELEAY; ISSN: 0040-4039
Elsevier Science B.V.
DOCUMENT TYPE:
Journal
LANGUAGE:
English
OTHER SOURCE(S):
CASREACT 139:337687

AB New gemini organogelators linked by an oxalyl amide that can be easily, effectively, and cheaply synthesized have good organogelation abilities and their cyclohexane gels have superior thermal stabilities; especially 7 possessing the branched alkyl ester can gel at 0.7 wt% cyclohexane even at 70°C.

IT 615584-80-0P 615584-81-1P 615584-82-2P
615584-86-6P
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(NMR and FT-IR on gelation of prepared gemini oxalyl-amide linked Inned
organogelators)
RN 615584-80-0 HCAPLUS
CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecy)- (9c1)
(CA INDEX NAME)

Absolute stereochemistry.

RN 615584-81-1 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, diethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-84-4 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, didodecyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-85-5 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(2-ethylhexyl) ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 615584-86-6 HCAPLUS CN L-Lysine, N2,N2'-(1,2-dioxo-1,2-ethanediyl)bis[N6-(1-oxododecyl)-, bis(3,5,5-trimethylhexyl) ester (9CI) (CA INDEX NAME) Absolute stereochemistry.

REFERENCE COUNT: AVAILABLE FOR THIS

36 THERE ARE 36 CITED REFERENCES

RECORD. ALL CITATIONS AVAILABLE IN THE

RE FORMAT

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(FILE 'HOME' ENTERED AT 15:17:16 ON 20 OCT 2004)

L1 L2 L3

FILE 'HCAPLUS' ENTERED AT 15:18:34 ON 20 OCT 2004 4 S L3 L4

FILE 'STNGUIDE' ENTERED AT 15:19:09 ON 20 OCT 2004

FILE 'REGISTRY' ENTERED AT 15:21:19 ON 20 OCT 2004 STRUCTURE UPLOADED 7 S L5 FUL L5 L6

FILE 'HCAPLUS' ENTERED AT 15:22:22 ON 20 OCT 2004  $^{\rm 2}$  S  $^{\rm L6}$ L7

FILE 'STNGUIDE' ENTERED AT 15:23:49 ON 20 OCT 2004

FILE 'REGISTRY' ENTERED AT 15:27:12 ON 20 OCT 2004

FILE 'STNGUIDE' ENTERED AT 15:31:14 ON 20 OCT 2004

FILE 'REGISTRY' ENTERED AT 15:32:29 ON 20 OCT 2004 STRUCTURE UPLOADED 17 S L8 FUL

L8 L9

FILE 'HCAPLUS' ENTERED AT 15:33:19 ON 20 OCT 2004  $^{\rm 2}$  S L9 L10

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	ENTRY
SESSION CA SUBSCRIBER PRICE 5.60	-1.40

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